

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN100519\
 Data File : BN008031.D
 Acq On : 05 Oct 2019 18:15
 Operator : JU
 Sample : K5078-04RE
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EXT-024-SW068-092619RE

Manual Integrations
 APPROVED

mohammad
 10/7/2019 4:39:52 PM

Quant Time: Oct 07 06:56:26 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN091819.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Thu Oct 03 18:39:49 2019

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	4820	0.40	ng	-0.02
7) Naphthalene-d8	10.44	136	19262	0.40	ng	-0.03
13) Acenaphthene-d10	14.30	164	12067	0.40	ng	-0.03
19) Phenanthrene-d10	17.05	188	24232	0.40	ng	-0.03
27) Chrysene-d12	21.25	240	25660	0.40	ng	-0.02
34) Perylene-d12	23.47	264	28815	0.40	ng	-0.03

System Monitoring Compounds

4) 2-Fluorophenol	5.30	112	1556	0.09	ng	-0.02
5) Phenol-d6	6.87	99	1812	0.08	ng	0.00
8) Nitrobenzene-d5	8.82	82	1849	0.11	ng	-0.01
11) 2-Methylnaphthalene-d10	12.04	152	3377	0.10	ng	-0.02
14) 2,4,6-Tribromophenol	15.80	330	647	0.10	ng	-0.02
15) 2-Fluorobiphenyl	12.93	172	4489	0.09	ng	-0.02
25) Fluoranthene-d10	19.09	212	7623	0.09	ng	-0.02
29) Terphenyl-d14	19.69	244	6471	0.10	ng	-0.02

Target Compounds

						Qvalue
9) Naphthalene	10.49	128	1692	0.031	ng	# 68
12) 2-Methylnaphthalene	12.11	142	1425	0.039	ng	# 83
16) Acenaphthylene	14.02	152	30433	0.476	ng	99
17) Acenaphthene	14.36	154	1446	0.035	ng	96
18) Fluorene	15.36	166	4345	0.082	ng	# 12
23) Phenanthrene	17.09	178	61638	0.817	ng	100
24) Anthracene	17.18	178	24810	0.364	ng	99
26) Fluoranthene	19.12	202	137550	1.465	ng	99
28) Pyrene	19.48	202	148043	1.692	ng	96
30) Benzo(a)anthracene	21.23	228	74972	0.793	ng	# 92
31) Chrysene	21.28	228	92697	1.007	ng	# 94
32) Bis(2-ethylhexyl)phthalate	21.18	149	40300	0.829	ng	98
33) Indeno(1,2,3-cd)pyrene	25.68	276	69128	0.600	ng	# 95
35) Benzo(b)fluoranthene	22.81	252	124426	1.244	ng	# 87
36) Benzo(k)fluoranthene	22.84	252	38237m	0.387	ng	
37) Benzo(a)pyrene	23.37	252	83926	0.893	ng	# 88
38) Dibenzo(a,h)anthracene	25.68	278	20773	0.216	ng	# 45
39) Benzo(g,h,i)perylene	26.36	276	83732	0.880	ng	93

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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