

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE032319\
 Data File : BE099129.D
 Acq On : 24 Mar 2019 2:30
 Operator : JU/SJ
 Sample : K1948-04
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PST-026-B020-031319

Manual Integrations
 APPROVED

Sohil
 3/25/2019 12:37:27 PM

Quant Time: Mar 24 03:55:16 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE032219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 23 04:40:11 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	4038	0.40	ng	0.00
7) Naphthalene-d8	10.63	136	15716	0.40	ng	-0.01
13) Acenaphthene-d10	14.49	164	10665	0.40	ng	0.00
19) Phenanthrene-d10	17.21	188	32131	0.40	ng	0.00
27) Chrysene-d12	21.38	240	42690	0.40	ng	-0.01
34) Perylene-d12	23.89	264	45910	0.40	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Fluorophenol	5.41	112	4913	0.41	ng	0.00
5) Phenol-d6	6.99	99	7138	0.41	ng	0.00
8) Nitrobenzene-d5	8.99	82	5225	0.76	ng	0.00
11) 2-Methylnaphthalene-d10	12.22	152	11210	0.40	ng	-0.01
14) 2,4,6-Tribromophenol	15.96	330	2405	0.77	ng	0.00
15) 2-Fluorobiphenyl	13.10	172	15827	0.37	ng	-0.01
25) Fluoranthene-d10	19.24	212	171164	0.40	ng	0.00
29) Terphenyl-d14	19.83	244	29216	0.32	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
16) Acenaphthylene	14.20	152	2161	0.038	ng	# 94
23) Phenanthrene	17.25	178	18490	0.194	ng	98
24) Anthracene	17.35	178	2086	0.023	ng	# 94
26) Fluoranthene	19.27	202	51246	0.382	ng	98
28) Pyrene	19.63	202	52661	0.364	ng	95
30) Benzo(a)anthracene	21.37	228	27722	0.184	ng	96
31) Chrysene	21.41	228	31155	0.207	ng	94
32) Bis(2-ethylhexyl)phthalate	21.29	149	46742	0.239	ng	99
33) Indeno(1,2,3-cd)pyrene	26.54	276	18709	0.116	ng	98
35) Benzo(b)fluoranthene	23.12	252	39509	0.249	ng	# 93
36) Benzo(k)fluoranthene	23.17	252	13870m	0.089	ng	
37) Benzo(a)pyrene	23.78	252	26047	0.177	ng	97
38) Dibenzo(a,h)anthracene	26.57	278	4659	0.033	ng	# 62
39) Benzo(g,h,i)perylene	27.37	276	19321	0.132	ng	98

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

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