

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN081219\
 Data File : BN007119.D
 Acq On : 13 Aug 2019 07:31
 Operator : HP/JU
 Sample : K4240-01
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :

Quant Time: Aug 13 08:48:43 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN081019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Aug 10 21:44:50 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.43	152	9459	0.40	ng	0.00
6) Naphthalene-d8	10.19	136	36749	0.40	ng	0.00
12) Acenaphthene-d10	14.08	164	20967	0.40	ng	0.00
18) Phenanthrene-d10	16.82	188	49444	0.40	ng	0.00
26) Chrysene-d12	21.05	240	48110	0.40	ng	-0.01
32) Perylene-d12	23.16	264	53127	0.40	ng	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 2-Fluorophenol	5.06	112	5251	0.23	ng	0.00
4) Phenol-d6	6.62	99	7176	0.25	ng	0.00
7) Nitrobenzene-d5	8.56	82	5728	0.19	ng	0.00
10) 2-Methylnaphthalene-d10	11.79	152	12446	0.18	ng	0.00
13) 2,4,6-Tribromophenol	15.57	330	1848	0.16	ng	0.00
14) 2-Fluorobiphenyl	12.71	172	3198	0.10	ng	0.00
24) Fluoranthene-d10	18.87	212	27884	0.16	ng	0.00
28) Terphenyl-d14	19.49	244	21253	0.16	ng	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
15) Acenaphthylene	13.79	152	5751	0.052	ng	100
22) Phenanthrene	16.87	178	17796	0.120	ng	98
23) Anthracene	16.95	178	4064	0.030	ng	# 91
25) Fluoranthene	18.90	202	52008	0.274	ng	99
27) Pyrene	19.26	202	49402	0.293	ng	# 96
29) Benzo(a)anthracene	21.02	228	30099	0.171	ng	# 88
30) Chrysene	21.08	228	33138	0.194	ng	# 91
31) Indeno(1,2,3-cd)pyrene	25.24	276	23080	0.123	ng	99
33) Benzo(b)fluoranthene	22.54	252	45346	0.249	ng	# 80
34) Benzo(k)fluoranthene	22.58	252	14367m	0.080	ng	
35) Benzo(a)pyrene	23.07	252	30259	0.183	ng	# 79
36) Dibenzo(a,h)anthracene	25.25	278	6166	0.037	ng	# 4
37) Benzo(a,h,i)perylene	25.87	276	23233	0.137	ng	# 83

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

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