

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE071718\
 Data File : BE096960.D
 Acq On : 17 Jul 2018 16:06
 Operator : SJ/JU
 Sample : J3986-05
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Jul 18 03:15:25 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE071718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 18 03:13:42 2018
 Response via : Initial Calibration

Internal Standards	R.T.	OIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	1559	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	7130	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	3726	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	10977	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	12329	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	11996	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	3896	0.34	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	12023	0.36	ng/ul	0.00

Target Compounds	R.T.	OIon	Response	Conc	Units	Ovalue
3) Naphthalene	10.19	128	5272	0.305	ng/ul#	93
5) 2-Methylnaphthalene	11.81	142	6310	0.553	ng/ul	98
7) Acenaphthylene	13.74	152	4133	0.254	ng/ul	98
8) Acenaphthene	14.09	153	41	0.003	ng/ul#	71
9) Fluorene	15.08	166	6777	0.506	ng/ul	91
11) Pentachlorophenol	16.43	266	914	0.737	ng/ul	97
12) Phenanthrene	16.82	178	8790	0.305	ng/ul#	89
13) Anthracene	16.91	178	12558	0.499	ng/ul#	93
15) Fluoranthene	18.84	202	10026	0.264	ng/ul	84
17) Pyrene	19.21	202	13453	0.377	ng/ul#	81
18) Benzo(a)anthracene	20.96	228	14730	0.451	ng/ul#	85
19) Chrysene	21.02	228	16845	0.455	ng/ul	99
21) Benzo(b)fluoranthene	22.54	252	22854	0.671	ng/ul	83
22) Benzo(k)fluoranthene	22.58	252	11427	0.325	ng/ul#	85
23) Benzo(a)pyrene	23.11	252	8910	0.298	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	25.53	276	5453	0.163	ng/ul#	79
25) Dibenzo(a,h)anthracene	25.53	278	21314	0.762	ng/ul#	87
26) Benzo(a,h,i)perylene	26.20	276	9664	0.323	ng/ul	96

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

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