

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE101018\
 Data File : BE097822.D
 Acq On : 10 Oct 2018 10:59
 Operator : SJ/JU
 Sample : SST00.838
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST00.838

Quant Time: Oct 10 11:31:14 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE101018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 10 11:08:35 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	2860	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	12549	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	7439	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	20704	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	36428	0.40	ng/ul	0.00
20) Perylene-d12	24.03	264	27443	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	16499	0.85	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	57091	0.84	ng/ul	0.00

Target Compounds						Qvalue
3) Naphthalene	10.74	128	24904	0.814	ng/ul	100
5) 2-Methylnaphthalene	12.36	142	17081	0.883	ng/ul	99
7) Acenaphthylene	14.26	152	21469	0.792	ng/ul	99
8) Acenaphthene	14.60	153	18131	0.778	ng/ul	98
9) Fluorene	15.59	166	22070	0.827	ng/ul	99
11) Pentachlorophenol	16.94	266	1421	0.697	ng/ul	98
12) Phenanthrene	17.34	178	41782	0.784	ng/ul	99
13) Anthracene	17.43	178	36472	0.814	ng/ul	97
15) Fluoranthene	19.36	202	56400	0.824	ng/ul	99
17) Pyrene	19.72	202	68668	0.700	ng/ul	99
18) Benzo(a)anthracene	21.46	228	76918	0.822	ng/ul	99
19) Chrysene	21.52	228	72544	0.673	ng/ul	98
21) Benzo(b)fluoranthene	23.25	252	59975	0.816	ng/ul	97
22) Benzo(k)fluoranthene	23.29	252	70045	0.855	ng/ul#	97
23) Benzo(a)pyrene	23.91	252	55800	0.821	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	26.72	276	64595	0.733	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.74	278	54795	0.734	ng/ul	99
26) Benzo(g,h,i)perylene	27.53	276	59379	0.790	ng/ul	99

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

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