

Data Path : Z:\HPCHEM1\BNA_M\Data\BM092916\
 Data File : BM007624.D
 Acq On : 29 Sep 2016 13:23
 Operator : UM/SJ
 Sample : H4955-21
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4VM9

Quant Time: Sep 29 14:54:59 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM092116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 23 04:00:23 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	1301	0.40	ng/ul	0.00
2) Naphthalene-d8	10.57	136	4839	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.42	164	3546	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.16	188	7957	0.40	ng/ul	0.00
16) Chrysene-d12	21.35	240	7229	0.40	ng/ul	0.00
20) Perylene-d12	23.62	264	5291	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.19	152	2046	0.36	ng/ul	0.02
14) Fluoranthene-d10	19.19	212	6986	0.38	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.62	128	10186	0.90	ng/ul	98
7) Acenaphthylene	14.13	152	10268	0.68	ng/ul	95
8) Acenaphthene	14.47	153	8284	0.71	ng/ul	89
9) Fluorene	15.48	166	12321	0.86	ng/ul	98
11) Pentachlorophenol	16.82	266	4647	2.14	ng/ul	97
12) Phenanthrene	17.19	178	16537	0.77	ng/ul	96
13) Anthracene	17.30	178	9099	0.47	ng/ul	97
17) Pyrene	19.58	202	41969	1.76	ng/ul	99
18) Benzo(a)anthracene	21.32	228	14389	0.67	ng/ul	97
19) Chrysene	21.38	228	24005	1.05	ng/ul	95
21) Benzo(b)fluoranthene	22.97	252	9199	0.52	ng/ul	95
22) Benzo(k)fluoranthene	22.97	252	9310	0.52	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.91	276	12352	0.73	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.91	278	518	0.04	ng/ul#	1
26) Benzo(g,h,i)perylene	26.60	276	14306	0.94	ng/ul	97

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

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