

Data Path : Z:\HPCHEM1\BNA M\DATA\BM110716\
 Data File : BM007776.D
 Acq On : 07 Nov 2016 17:58
 Operator : UM/SJ
 Sample : H5552-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AA2

Manual Integrations
 APPROVED

umangi
 11/8/2016 3:15:21 PM

Quant Time: Nov 08 03:06:05 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM102016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 08 02:10:44 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	173198	20.00	ng/ul	0.00
18) Naphthalene-d8	10.53	136	669601	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.38	164	435353	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.13	188	881772	20.00	ng/ul	0.00
75) Chrysene-d12	21.31	240	1132317	20.00	ng/ul	0.00
83) Perylene-d12	23.56	264	1175646	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.29	96	5801	1.44	ng/uL	0.00
5) Phenol-d5	6.92	99	91767	6.86	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.09	67	251475	31.70	ng/ul	0.00
9) 2-Chlorophenol-d4	7.28	132	275823	26.90	ng/ul	0.00
13) 4-Methylphenol-d8	8.45	113	175579	16.79	ng/ul	0.00
19) Nitrobenzene-d5	8.90	128	164315	34.10	ng/ul	0.00
22) 2-Nitrophenol-d4	9.62	143	181002	35.43	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.15	165	305950	30.46	ng/ul	0.00
29) 4-Chloroaniline-d4	10.71	131	4005m	0.35	ng/ul	0.04
43) Dimethylphthalate-d6	13.79	166	1045560	35.27	ng/ul	0.00
46) Acenaphthylene-d8	14.07	160	1240235	34.76	ng/ul	0.00
51) 4-Nitrophenol-d4	14.62	143	27697m	5.91	ng/ul	0.02
57) Fluorene-d10	15.37	176	904756	34.82	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.51	200	146652	27.99	ng/ul	0.00
70) Anthracene-d10	17.22	188	1411784	35.53	ng/ul	0.00
76) Pyrene-d10	19.52	212	1660675	35.70	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.42	264	1805173	35.17	ng/ul	0.00

Target Compounds	Ovalue

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

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