

Data Path : Z:\HPCHEM1\BNA M\DATA\BM121716\
 Data File : BM008376.D
 Acq On : 16 Dec 2016 16:48
 Operator : UM/SJ
 Sample : H6039-03
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA8R9

Manual Integrations
 APPROVED

umangi
 12/20/2016 10:42:00 AM

Quant Time: Dec 17 02:21:51 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-2016\SOM02.2-EPA-BM121
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Dec 17 02:12:55 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	104796	20.00	ng/ul	0.00
18) Naphthalene-d8	10.44	136	414681	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.30	164	274018	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.06	188	551347	20.00	ng/ul	0.00
75) Chrysene-d12	21.26	240	683501	20.00	ng/ul	0.00
83) Perylene-d12	23.49	264	702893	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.23	96	4566	2.05	ng/uL	0.00
5) Phenol-d5	6.86	99	73401	9.03	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	7.01	67	184469	39.48	ng/ul	0.00
9) 2-Chlorophenol-d4	7.20	132	207996	34.00	ng/ul	0.00
13) 4-Methylphenol-d8	8.38	113	133766	21.24	ng/ul	0.00
19) Nitrobenzene-d5	8.82	128	126421	42.19	ng/ul	0.00
22) 2-Nitrophenol-d4	9.54	143	140780	42.39	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.09	165	227059	37.31	ng/ul	0.01
29) 4-Chloroaniline-d4	10.69	131	2256m	0.32	ng/ul	0.09
43) Dimethylphthalate-d6	13.72	166	781404	42.90	ng/ul	0.00
46) Acenaphthylene-d8	13.99	160	970172	45.07	ng/ul	0.00
51) 4-Nitrophenol-d4	14.63	143	14971m	5.58	ng/ul	0.07
57) Fluorene-d10	15.30	176	676755	42.74	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.47	200	96237	29.32	ng/ul	0.00
70) Anthracene-d10	17.16	188	1055865	43.96	ng/ul	0.00
76) Pyrene-d10	19.46	212	1225929	46.44	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.34	264	1250893	42.38	ng/ul	0.00

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

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