

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM022316\
 Data File : BM004357.D
 Acq On : 23 Feb 2016 19:34
 Operator : SJ/UM
 Sample : H1542-01
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 0311-RIN05

Manual Integrations
 APPROVED

UMANGI
 2/24/2016 4:31:07 PM

Quant Time: Feb 24 02:48:46 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM022216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Feb 24 01:02:10 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	28458	20.00	ng/ul	0.00
18) Naphthalene-d8	10.41	136	134232	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.29	164	85019	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.04	188	200449	20.00	ng/ul	0.00
78) Chrysene-d12	21.25	240	197155	20.00	ng/ul	0.00
86) Perylene-d12	23.47	264	146261	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.09	96	598	1.11	ng/uL	0.00
5) Phenol-d5	6.82	99	15128m	6.34	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.96	67	40772	33.60	ng/ul	0.00
9) 2-Chlorophenol-d4	7.16	132	53692	26.09	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	33111	15.53	ng/ul	0.00
19) Nitrobenzene-d5	8.78	128	34033	33.33	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	39310	33.77	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.05	165	61252m	29.11	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0	0.00	ng/ul	
44) Dimethylphthalate-d6	13.70	166	260926	35.79	ng/ul	0.00
47) Acenaphthylene-d8	13.97	160	304385	35.66	ng/ul	0.00
52) 4-Nitrophenol-d4	14.55	143	4249m	3.63	ng/ul	0.02
58) Fluorene-d10	15.28	176	224738	36.24	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.43	200	35353m	28.71	ng/ul	0.00
71) Anthracene-d10	17.14	188	367468	37.41	ng/ul	0.00
79) Pyrene-d10	19.44	212	404066	42.37	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.33	264	286621	37.34	ng/ul	0.00

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

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