

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111419\
 Data File : BM023729.D
 Acq On : 14 Nov 2019 22:41
 Operator : JU
 Sample : K5700-07
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BEGZ9

Quant Time: Nov 15 06:34:07 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM11119MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 15 06:30:14 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.53	152	329870	20.00	ng/ul	0.00
18) Naphthalene-d8	10.30	136	1348312	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.19	164	873219	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.94	188	1895685	20.00	ng/ul	0.00
78) Chrysene-d12	21.14	240	1858984	20.00	ng/ul	0.00
86) Perylene-d12	23.30	264	2184728	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.06	96	51394	6.42	ng/uL	0.00
5) Phenol-d5	6.72	99	151286	5.03	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.87	67	433784	25.11	ng/ul	0.00
9) 2-Chlorophenol-d4	7.07	132	472080	21.67	ng/ul	0.00
13) 4-Methylphenol-d8	8.25	113	289544	12.07	ng/ul	0.00
19) Nitrobenzene-d5	8.68	128	295870	30.66	ng/ul	0.00
22) 2-Nitrophenol-d4	9.40	143	338685	33.39	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.94	165	592408	25.61	ng/ul	0.00
29) 4-Chloroaniline-d4	10.46	131	36514	1.44	ng/ul	0.00
44) Dimethylphthalate-d6	13.61	166	2067934	30.37	ng/ul	0.00
47) Acenaphthylene-d8	13.87	160	2366934	30.46	ng/ul	0.00
52) 4-Nitrophenol-d4	14.42	143	47484	4.08	ng/ul	0.00
58) Fluorene-d10	15.19	176	1811269	30.12	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.32	200	270664	24.68	ng/ul	0.00
71) Anthracene-d10	17.03	188	2897848	32.22	ng/ul	0.00
79) Pyrene-d10	19.34	212	3199875	33.01	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.16	264	3708034	32.25	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
45) Dimethylphthalate	13.66	163	109055	1.634	ng/ul	99

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

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