

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
 Data File : BM025804.D
 Acq On : 26 Mar 2020 17:39
 Operator : CG/JU
 Sample : PB127667BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK67

Quant Time: Mar 26 18:35:05 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 24 00:45:06 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	218993	20.00	ng/ul	-0.01
18) Naphthalene-d8	10.33	136	898227	20.00	ng/ul	-0.02
36) Acenaphthene-d10	14.20	164	549118	20.00	ng/ul	-0.01
62) Phenanthrene-d10	16.95	188	1184957	20.00	ng/ul	-0.01
78) Chrysene-d12	21.14	240	1305307	20.00	ng/ul	-0.01
86) Perylene-d12	23.30	264	1474880	20.00	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.13	96	38202	7.32	ng/uL	0.00
5) Phenol-d5	6.75	99	616101	35.65	ng/ul	-0.01
7) Bis-(2-Chloroethyl)ether-d	6.91	67	365873	37.09	ng/ul	-0.01
9) 2-Chlorophenol-d4	7.10	132	523718	37.34	ng/ul	-0.01
13) 4-Methylphenol-d8	8.28	113	507661	35.75	ng/ul	-0.01
19) Nitrobenzene-d5	8.71	128	262332	39.29	ng/ul	0.00
22) 2-Nitrophenol-d4	9.42	143	278854	39.29	ng/ul	-0.01
26) 2,4-Dichlorophenol-d3	9.96	165	519537	35.80	ng/ul	-0.01
29) 4-Chloroaniline-d4	10.47	131	740613	43.94	ng/ul	-0.01
44) Dimethylphthalate-d6	13.62	166	1619261	38.80	ng/ul	-0.01
47) Acenaphthylene-d8	13.89	160	1963094	39.26	ng/ul	-0.01
52) 4-Nitrophenol-d4	14.41	143	270808	34.08	ng/ul	-0.01
58) Fluorene-d10	15.20	176	1426856	38.95	ng/ul	-0.01
63) 4,6-Dinitro-2-methylphenol	15.33	200	219670	30.38	ng/ul	-0.01
71) Anthracene-d10	17.05	188	2220205	40.11	ng/ul	0.00
79) Pyrene-d10	19.35	212	2430197	38.76	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.16	264	3032956	40.17	ng/ul	-0.01

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

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