

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032320\
 Data File : BM025799.D
 Acq On : 26 Mar 2020 14:32
 Operator : CG/JU
 Sample : PB127634BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK34

Quant Time: Mar 26 15:22:34 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	7.57	152	173141	20.00	ng/ul	0.00
18) Naphthalene-d8	10.33	136	697485	20.00	ng/ul	-0.01
36) Acenaphthene-d10	14.20	164	427649	20.00	ng/ul	-0.01
62) Phenanthrene-d10	16.95	188	926744	20.00	ng/ul	-0.01
78) Chrysene-d12	21.14	240	1023213	20.00	ng/ul	-0.01
86) Perylene-d12	23.29	264	1158484	20.00	ng/ul	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.14	96	28803	6.98	ng/uL	0.00
5) Phenol-d5	6.75	99	487663	35.69	ng/ul	-0.01
7) Bis-(2-Chloroethyl)ether-d	6.91	67	293768	37.66	ng/ul	-0.01
9) 2-Chlorophenol-d4	7.10	132	418983	37.78	ng/ul	-0.01
13) 4-Methylphenol-d8	8.27	113	402402	35.84	ng/ul	-0.01
19) Nitrobenzene-d5	8.71	128	206019	39.74	ng/ul	0.00
22) 2-Nitrophenol-d4	9.43	143	216585	39.30	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.96	165	404197	35.86	ng/ul	-0.01
29) 4-Chloroaniline-d4	10.47	131	582632	44.52	ng/ul	-0.01
44) Dimethylphthalate-d6	13.63	166	1276670	39.28	ng/ul	0.00
47) Acenaphthylene-d8	13.89	160	1539839	39.54	ng/ul	-0.01
52) 4-Nitrophenol-d4	14.41	143	204673	33.08	ng/ul	-0.01
58) Fluorene-d10	15.20	176	1123602	39.38	ng/ul	-0.01
63) 4,6-Dinitro-2-methylphenol	15.33	200	164356	29.07	ng/ul	-0.01
71) Anthracene-d10	17.05	188	1744580	40.30	ng/ul	0.00
79) Pyrene-d10	19.35	212	1915087	38.97	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.16	264	2417226	40.75	ng/ul	-0.01

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

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