

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP030321\
 Data File : BP004892.D
 Acq On : 03 Mar 2021 14:19
 Operator : CG/JU
 Sample : PB134916BL
 Misc : SOM-MDL-S-BL01
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK16

Quant Time: Mar 03 14:50:51 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP030221MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Mar 03 10:35:24 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	28810	20.00	ng/ul	0.00
18) Naphthalene-d8	10.53	136	111966	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.39	164	78402	20.00	ng/ul	-0.01
62) Phenanthrene-d10	17.15	188	183458	20.00	ng/ul	-0.01
78) Chrysene-d12	21.24	240	200857	20.00	ng/ul	-0.01
86) Perylene-d12	23.54	264	207739	20.00	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.24	96	4627	6.32	ng/uL	0.00
5) Phenol-d5	6.92	99	75321	32.10	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.08	67	41624	34.21	ng/ul	-0.01
9) 2-Chlorophenol-d4	7.28	132	60579	33.47	ng/ul	0.00
13) 4-Methylphenol-d8	8.45	113	60597	32.41	ng/ul	-0.01
19) Nitrobenzene-d5	8.90	128	28832	35.35	ng/ul	0.00
22) 2-Nitrophenol-d4	9.62	143	32087	35.04	ng/ul	-0.01
26) 2,4-Dichlorophenol-d3	10.16	165	57720	33.33	ng/ul	-0.01
29) 4-Chloroaniline-d4	10.67	131	80277	40.55	ng/ul	0.00
44) Dimethylphthalate-d6	13.80	166	202830	35.36	ng/ul	-0.01
47) Acenaphthylene-d8	14.09	160	251729	35.85	ng/ul	0.00
52) 4-Nitrophenol-d4	14.60	143	32602	31.77	ng/ul	-0.01
58) Fluorene-d10	15.39	176	189138	38.78	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.51	200	38478	32.17	ng/ul	-0.01
71) Anthracene-d10	17.25	188	313554	37.69	ng/ul	0.00
79) Pyrene-d10	19.49	212	371139	39.15	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.39	264	423145	39.74	ng/ul	-0.01

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

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