

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP120820\
 Data File : BP004192.D
 Acq On : 08 Dec 2020 09:52
 Operator : CG/JU
 Sample : PB133372BL
 Misc : MDL-SOIL-BL-06
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK372

Quant Time: Dec 08 10:57:14 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP120420.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 08 06:09:39 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	378457	20.00	ng/ul	0.00
20) Naphthalene-d8	10.64	136	1567374	20.00	ng/ul	-0.01
38) Acenaphthene-d10	14.49	164	965173	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.25	188	2060209	20.00	ng/ul	0.00
79) Chrysene-d12	21.34	240	2074241	20.00	ng/ul	-0.01
88) Perylene-d12	23.70	264	2211620	20.00	ng/ul	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.34	96	79692	8.09	ng/uL	0.00
4) Pyridine-d5	3.73	84	1097537	39.22	ng/ul	-0.01
7) Phenol-d5	7.00	99	1257031	37.15	ng/ul	-0.01
9) Bis-(2-Chloroethyl)ether-d	7.18	67	792394	40.51	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.38	132	970581	36.86	ng/ul	-0.01
15) 4-Methylphenol-d8	8.54	113	969814	36.38	ng/ul	-0.01
21) Nitrobenzene-d5	9.00	128	514356	40.25	ng/ul	0.00
24) 2-Nitrophenol-d4	9.72	143	435598	37.33	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.25	165	924342	36.00	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.77	131	1424433	37.81	ng/ul	-0.01
46) Dimethylphthalate-d6	13.90	166	3147498	41.06	ng/ul	-0.01
49) Acenaphthylene-d8	14.18	160	3688774	39.16	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.67	143	430545	34.32	ng/ul	0.00
60) Fluorene-d10	15.49	176	2681278	40.00	ng/ul	-0.01
65) 4,6-Dinitro-2-methylphenol	15.60	200	272709	32.58	ng/ul	-0.01
73) Anthracene-d10	17.35	188	4193177	41.41	ng/ul	-0.01
81) Pyrene-d10	19.59	212	4728202	43.33	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.56	264	4981919	41.93	ng/ul	-0.01

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

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