

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090921\
 Data File : BP006944.D
 Acq On : 10 Sep 2021 16:54
 Operator : CG/JU
 Sample : PB138976BL
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK976

Quant Time: Sep 10 17:36:09 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP090821.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 09 11:07:24 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	316439	20.000	ng/u1	0.00
20) Naphthalene-d8	10.734	136	1279116	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.557	164	785295	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.310	188	1620787	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.392	240	1557145	20.000	ng/u1	# 0.00
88) Perylene-d12	23.821	264	1571760	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	46245	5.816	ng/uL	0.00
4) Pyridine-d5	3.787	84	612927	29.579	ng/u1	0.00
7) Phenol-d5	7.093	99	710778	30.561	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	424989	31.817	ng/u1	0.00
11) 2-Chlorophenol-d4	7.463	132	578550	30.814	ng/u1	0.00
15) 4-Methylphenol-d8	8.634	113	549455	30.084	ng/u1	0.00
21) Nitrobenzene-d5	9.093	128	254020	32.786	ng/u1	0.00
24) 2-Nitrophenol-d4	9.816	143	246133	32.268	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.352	165	514348	28.597	ng/u1	0.00
31) 4-Chloroaniline-d4	10.869	131	761392	30.612	ng/u1	0.00
46) Dimethylphthalate-d6	13.969	166	1654112	32.522	ng/u1	0.00
49) Acenaphthylene-d8	14.251	160	1966771	30.532	ng/u1	0.00
54) 4-Nitrophenol-d4	14.740	143	233762	26.455	ng/u1	0.00
60) Fluorene-d10	15.551	176	1417100	31.569	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	154904	21.605	ng/u1	0.00
73) Anthracene-d10	17.404	188	2215300	33.014	ng/u1	-0.01
81) Pyrene-d10	19.639	212	2548438	34.946	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.668	264	2401167	31.667	ng/u1	-0.01

Target Compounds	Qvalue

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

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