

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072621\
 Data File : BP006347.D
 Acq On : 26 Jul 2021 20:17
 Operator : CG/JU
 Sample : PB137859BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK859

Quant Time: Jul 27 01:10:00 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jul 24 04:29:12 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	278775	20.000	ng/u1	0.00
20) Naphthalene-d8	10.569	136	1233881	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.410	164	723259	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.163	188	1419214	20.000	ng/u1	0.00
79) Chrysene-d12	21.263	240	1311443	20.000	ng/u1	0.00
88) Perylene-d12	23.610	264	1342499	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	56670	7.588	ng/uL	0.00
4) Pyridine-d5	3.699	84	857630	38.683	ng/u1	0.00
7) Phenol-d5	6.958	99	980842	37.462	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	644915	39.588	ng/u1	0.00
11) 2-Chlorophenol-d4	7.322	132	761070	39.126	ng/u1	0.00
15) 4-Methylphenol-d8	8.493	113	783784	38.067	ng/u1	0.00
21) Nitrobenzene-d5	8.940	128	372826	40.580	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.657	143	376886	42.963	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.193	165	636043	37.000	ng/u1	0.00
31) 4-Chloroaniline-d4	10.710	131	1083523	38.286	ng/u1	0.00
46) Dimethylphthalate-d6	13.834	166	2186916	42.484	ng/u1	0.00
49) Acenaphthylene-d8	14.104	160	2594263	41.021	ng/u1	0.00
54) 4-Nitrophenol-d4	14.610	143	400745	39.208	ng/u1	0.00
60) Fluorene-d10	15.410	176	1754061	40.534	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.528	200	263653	33.324	ng/u1	0.00
73) Anthracene-d10	17.263	188	2748742	42.905	ng/u1	0.00
81) Pyrene-d10	19.504	212	2946028	43.675	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.457	264	2814060	41.249	ng/u1	0.00

Target Compounds Qvalue

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

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